

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA E\DATA\BE101417\
 Data File : BE094377.D
 Acq On : 15 Oct 2017 4:34
 Operator : SJ/JU
 Sample : I5548-21
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AE7

Manual Integrations
 APPROVED

Sohil
 10/16/2017 8:25:42 PM

Quant Time: Oct 16 05:52:25 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 16 05:09:03 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1869	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	9878	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	5989	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	12172	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	11492	0.40	ng/ul	0.00
20) Perylene-d12	23.95	264	11740m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	3600	0.24	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	8189	0.20	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.74	128	80941	3.345	ng/ul#	91
5) 2-Methylnaphthalene	12.35	142	27676	1.763	ng/ul	98
7) Acenaphthylene	14.24	152	5959	0.255	ng/ul#	93
8) Acenaphthene	14.58	153	113099	5.884	ng/ul	97
9) Fluorene	15.57	166	113948	5.062	ng/ul#	92
12) Phenanthrene	17.29	178	923641	29.038	ng/ul#	85
13) Anthracene	17.39	178	267381	8.643	ng/ul#	91
15) Fluoranthene	19.31	202	1136165	27.185	ng/ul	84
17) Pyrene	19.68	202	883360	27.627	ng/ul#	81
18) Benzo(a)anthracene	21.40	228	547882	17.971	ng/ul#	86
19) Chrysene	21.46	228	462593	13.139	ng/ul	99
21) Benzo(b)fluoranthene	23.18	252	491079	15.398	ng/ul	85
22) Benzo(k)fluoranthene	23.22	252	201915m	5.477	ng/ul	
23) Benzo(a)pyrene	23.84	252	376847m	11.371	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.63	276	237244m	6.998	ng/ul	
25) Dibenzo(a,h)anthracene	26.63	278	71868m	2.527	ng/ul	
26) Benzo(a,h,i)perylene	27.47	276	178465m	6.103	ng/ul	

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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